



Chemical and Environmental Measurement Information

Recra LabNet Philadelphia Analytical Report **REVISION**

Client: TNU-HANFORD B99-085

RFW#: 9909L126

SDG/SAF #: H0535/B99-085

W.O. #: 10985-001-001-9999-00

Date Received: 09-17-99



EDMC

SEMIVOLATILE

This narrative was corrected to add the TIC search for Tributylphosphate.

One (1) water sample was collected on 09-15-99.

The sample and its associated QC samples were extracted on 09-21-99 and analyzed according to criteria set forth in Recra OPs based on SW 846 Method 8270B TCL Semivolatile target compounds on 10-04-99.

The following is a summary of the QC results accompanying the sample results and a description of any problems encountered during their analyses:

1. The cooler temperature upon receipt has been recorded on the chain-of-custody.

- 2. The required holding times for extraction and analysis were met.
- 3. Non-target compounds were detected in these samples.
- 4. These samples were spectrally searched for Butylated Hydroxytoluene and Tributylphosphate; however, they were not identified in the samples.
- 5. All surrogate recoveries were within USEPA QC limits.
- 6. Two (2) of eleven (11) matrix spike recoveries were outside USEPA QC limits. A copy of the Sample Discrepancy Report (SDR) has been enclosed.
- 7. Two (2) of eleven (11) blank spike recoveries were outside USEPA QC limits. A copy of the Sample Discrepancy Report (SDR) has been enclosed.

J. Michael Taylor

Vice President

Philadelphia Analytical Laboratory

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0)-22-00 Date

The results presented in this report relate only to the analytical testing and conditions of the samples at receipt and during storage. All pages of this report are integral parts of the analytical data. Therefore, this report should only be reproduced in its entirety of 11 pages.

GLOSSARY OF BNA DATA

DATA QUALIFIERS

U	=	Compound was analyzed for but not detected. The associated numerical value is the estimated
		sample quantitation limit which is included and corrected for dilution and percent moisture.

- J = Indicates an estimated value. This flag is used under the following circumstances: 1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; or 2) when the mass spectral data indicate the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. For example, if the limit of detection is 10 ug/L and a concentration of 3 ug/L is calculated, it is reported as 3J.
- B = This flag is used when the analyte is found in the associated blank as well as in the sample. It indicates possible/probable blank contamination. This flag is also used for a TIC as well as for a positively identified TCL compound.
- E = Indicates that the compound was detected beyond the calibration range and was subsequently analyzed at a dilution.
- **D** = Identifies all compounds identified in an analysis at a secondary dilution factor.
- I = Interference.
- NQ = Result qualitatively confirmed but not able to quantify.
- A = Indicates that a TIC is a suspected aldol-condensation product.
- N = Indicates presumptive evidence of a compound. This flag is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It is applied to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the N code is not used.
- X = This flag is used for a TIC compound which is quantified relative to a response factor generated from a daily calibration standard (rather than quantified relative to the closest internal standard).
- Y = Additional qualifiers used as required are explained in the case narrative.

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GLOSSARY OF BNA DATA

ABBREVIATIONS

BS = Indicates blank spike in which reagent grade water is spiked with the CLP matrix spike solutions and carried through all the steps in the method. Spike recoveries are reported.

BSD = Indicates blank spike duplicate.

MS = Indicates matrix spike.

MSD = Indicates matrix spike duplicate.

DL = Suffix added to sample number to indicate that results are from a diluted analysis.

NA = Not Applicable.

DF = Dilution Factor.

NR = Not Required.

SP, Z = Indicates Spiked Compound.

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Initiator: Studie RFW Batch: 9909L124 Parameter: BNF Date: 10-5-99 Samples: MSD BS Matrix: Water Prep Batch: 991E1150 Conf
1. Reason for SDR a. COC Discrepancy Tech Profile Error Client Request Sampler Error on C-O-C Transcription Error Wrong Test Code Other b. General Discrepancy Missing Sample/Extract Container Broken Wrong Sample Pulled Label ID's Illegible Hold Time Exceeded Insufficient Sample Preservation Wrong Received Past Hold Improper Bottle Type Not Amenable to Analysis Note: Verified by [Log-In] or [Prep Group] (circle)signature/date: C. QC Problem (Include all relevant specific results; attach data if necessary) MSD & BS
2. Known or Probable Causes(s) possible problem with prep
3. Discussion and Proposed Action Re-log Entire Batch Following Samples: Re-leach Re-extract Re-digest Revise EDD Change Test Code to Place On/Take Off Hold (circle) 4. Project Manager Instructionssignature/date: Concur with Proposed Action Disagree with Proposed Action; See Instruction Include in Case Narrative Client Contacted: Date/Person Add Cancel
5. Final Actionsignature/date: (11) (0) 25/49 Other Explanation: Verified re-[log][leach][extract][digest][analysis] (circle) Included in Case Narrative Hard Copy COC Revised Electronic COC Revised EDD Corrections Completed When Final Action has been recorded, forward original to QA Specialist for distribution and filing.
Route Distribution of Completed SDR

Neura Labnet Philadelphia Sample Discrepancy Report (SDR) SDR#:

99 MS 081

Recra LabNet - Lionville Laboratory

Semivolatiles by GC/MS, HSL List

Client: TNU-HANFORD B99-085

RFW Batch Number: 9909L126

*= Outside of EPA CLP QC limits.

Report Date: 10/25/99 17:22

Page: la

Work Order: 10985001001

 \bigcirc Cust ID: BOWCP8 BOWCP8 BOWCP8 SBLKDI SBLKDI BS 99LE1150-MB1 Sample RFW#: 001 001 MS 001 MSD 99LE1150-MB1 Information WATER WATER WATER WATER WATER Matrix: DF. 1.00 1.00 1.00 1.00 1.00 Units: UG/L UG/L UG/L UG/L UG/L Nitrobenzene-d5 98 % કૃ % 2-Fluorobiphenyl 79 ۶ ٩. 75 ş 77 왓 67 è Surrogate 77 Recovery Terphenyl-d14 89 ş 93 왕 82 왕 82 87 % Phenol-d5 82 & 83 79 ş 69 81 2-Fluorophenol 70 å 79 2 75 ş 82 49 ş 2,4,6-Tribromophenol 73 ષ્ટ્ર 93 68 69 56 10 U 76 76 ٥. 10 II 65 Phenol bis(2-Chloroethyl)ether U 10 [] 10 U 20 U 20 10 U 2-Chlorophenol 10 U 73 કૃ 10 II 1,3-Dichlorobenzene 10 H 10 U 20 II 20 ΙŢ 10 II 1.4-Dichlorobenzene____ ş. 72 10 U 48 10 11 1,2-Dichlorobenzene 10 U 20 II 20 11 10 [] 10 U 2-Methylphenol_____ 10 [] 20 [] ΙŢ 10 II 10 U 2,2'-oxybis(1-Chloropropane) 10 U 20 U 20 10 U 10 U Ü 4-Methylphenol 10 U 20 U 20 U 10 U 10 U N-Nitroso-di-n-propylamine 10 U 98 ş 87 왕 10 U 71 Hexachloroethane_____ 10 U 20 U 20 U 10 U 10 U Nitrobenzene _____ 10 U 20 U 20 U 10 U 10 U Isophorone 10 U 20 U 20 U 10 U 10 II 2-Nitrophenol_____ 10 U 20 [] 20 II 10 II 10 U 2.4-Dimethylphenol____ 10 U 20 II 20 U 10 U 10 U bis (2-Chloroethoxy) methane 10 U 10 U 10 II 20 IJ 2.0 IJ 2,4-Dichlorophenol_____ 10 U 20 U 20 U 10 U 10 U 1,2,4-Trichlorobenzene____ 10 U 77 80 ջ 10 U 54 Naphthalene _____ 10 U Ħ 10 U 20 20 U 10 U 4-Chloroaniline _____ 10 U 10 U 10 IJ 20 U 20 U Hexachlorobutadiene____ 10 U 20 U 20 U 10 U 10 U 4-Chloro-3-methylphenol____ 10 U 10 U 72 81 73 2 2-Methylnaphthalene_____ 10 U 20 U 10 U 10 U 20 11 Hexachlorocyclopentadiene 20 U 10 II 20 U 10 U 10 U 2.4.6-Trichlorophenol_ 10 U 20 U 20 U 10 U 10 U 2.4.5-Trichlorophenol 25 U 50 U 50 U 25 U 25 U

RFW Batch Number: 9909L126	Client:	Client: TNU-HANFORD B99-085					k Order: 109	001001	Page: 1b	
Cust II	D: BOWCP	3	BOWCP8		BOWCP8		SBLKDI		SBLKDI BS	
RFW;	#: 00:	ı	001 MS		001 MSD		99LE1150-ME	31	99LE1150~MB	L
2-Chloronaphthalene	10	U U	20 t	 J	20	U	10	U	10 t	
2-Nitroaniline	25	U	50 t		50	U	25	U	25 U	
Dimethylphthalate	10		20 t		20	Ū	10	-	10 t	
Acenaphthylene	10		20 t		20	Ū	10		10 t	
2,6-Dinitrotoluene	10	U		- J	20	U	10	_	10 t	
3-Nitroaniline		Ū	50 T		50	Ū	25	_	25 (
Acenaphthene	10			} }	85	્ર	10	-	73 8	
2,4-Dinitrophenol	25	Ű		J	50	U	25	_	25 t	=
4-Nitrophenol		Ū		ì	0 *	_	25		8 * %	
Dibenzofuran	10	Ū	20 t		20	Ü	10	_	10 U	
2,4-Dinitrotoluene		Ū	99 * 8	-	84	%	10	-	69 %	
Diethylphthalate		Ū	20 t		20	Ü	10	-	10 t	
4-Chlorophenyl-phenylether	10	Ū	20 t		20	Ü	10		10 t	
Fluorene		U	20 T	-	20	Ü	10		10 t	-
-Nitroaniline	25	U		- J	50	Ū	25		25 t	-
,6-Dinitro-2-methylphenol	25	Ū	50 t	J	50	Ū	25		25 t	
N-Nitrosodiphenylamine (1)	10	Ū	20 t		20	Ū	10		10 U	J
-Bromophenyl-phenylether	10	U	20 t	J	20	Ū		Ū	10 t	
Mexachlorobenzene		U	20 t	J	20	Ū	10	Ū	10 U	J
Pentachlorophenol	25	U	73 %	<u>.</u>	17	%	25	Ü	7 * 8	•
Phenanthrene		U	20 t		20	Ü		Ū	10 0	J
Anthracene	10	Ū	20 t	J	20	Ū		Ū	10 U	
Carbazole		Ü		- J	20		10	_	10 0	
Di-n-butylphthalate		J		J	2	J	10		10 U	
Fluoranthene		Ū	20 t		20	Ū	10		10 U	
Pyrene		Ū	94 %		83	26	10	_	88 %	
Butylbenzylphthalate	10	Ū	20 L		20	Ū	10		10 U	
3,3'-Dichlorobenzidine	10	Ü	20 T	J	20	Ū	10	_	10 U	
Benzo(a)anthracene	10	Ü	20 T			Ū	10		10 U	
Chrysene		U	20 t		20		10	-	10 0	
ois(2-Ethylhexyl)phthalate		Ū	5 J			Ū	10	_	3 J	
i-n-octyl phthalate		Ū	20 U		20	_	10		10 U	
enzo(b)fluoranthene		Ū	20 U			Ü	10		10 U	
enzo(k)fluoranthene	10	Ū	20 U			Ū	10		10 U	
enzo(a)pyrene		Ü	20 U		20		10		10 U	
indeno(1,2,3-cd)pyrene		U	20 U		20		10		10 U	
Dibenz (a, h) anthracene		Ū	20 U			Ū		U	10 U	
Benzo(g,h,i)perylene			20 U			U		U	10 U	
(1) Cannot be generated from F			+ 0		20 21 D 20 1	, .		_	10 0	

Benzo(g,h,i)perylene_______ 10 U 20 U 20 U (1) - Cannot be separated from Diphenylamine. *= Outside of EPA CLP QC limits.

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CLIENT SAMPLE NO.

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

1		
BOWCP8		
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Lab Name: Recra.LabNet Work Order: 10985001001

Client: TNU-HANFORD B99-085

Lab Sample ID: 9909L126-001 Matrix: (soil/water) <u>WATER</u>

Sample wt/vol: $\underline{1000}$ (g/mL) $\underline{\text{ML}}$ Lab File ID: $\underline{\text{A100410}}$

Date Received: <u>09/17/99</u> Level: (low/med) LOW

% Moisture: ____ decanted: (Y/N)__ Date Extracted: 09/21/99

Concentrated Extract Volume: $1000 \,(\text{uL})$ Date Analyzed: 10/04/99

Dilution Factor: 1.00 Injection Volume: 2.0(uL)

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u> Number TICs found: 3

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	ĮΩĮ
		======	=======================================	====≠
1.	UNKNOWN	7.77	2	J
2.	UNKNOWN	7.94	3	J
3.	UNKNOWN	23.13	4	J
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CLIENT SAMPLE NO.

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

SBLKDI		
}		

Lab Name: Recra_LabNet Work Order: 10985001001

Client: TNU-HANFORD B99-085

Matrix: (soil/water) WATER Lab Sample ID: 99LE1150-MB1

Sample wt/vol: _1000 (g/mL) ML Lab File ID: A100408

Level: (low/med) LOW Date Received: 09/21/99

% Moisture: ____ decanted: (Y/N)__ Date Extracted: 09/21/99

Concentrated Extract Volume: 1000(uL) Date Analyzed: 10/04/99

Injection Volume: 2.0(uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: __7.0

CONCENTRATION UNITS: Number TICs found: 0 (ug/L or ug/Kg) <u>UG/L</u>

CAS NUME	BER	COMPOUND	NAME	RT	EST.	CONC.	Q
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Recra LabNet - Lionville Laboratory BNA ANALYTICAL DATA PACKAGE FOR TNU-HANFORD B99-085

RFW LOT # :9909L126

CLIENT ID RFW # MTX PREP # COLLECTION EXTR/PREP ANALYSIS

BOWCP8 001 W 99LE1150 09/15/99 09/21/99 10/04/99
BOWCP8 001 MS W 99LE1150 09/15/99 09/21/99 10/04/99
BOWCP8 001 MSD W 99LE1150 09/15/99 09/21/99 10/04/99

LAB QC:

DATE RECEIVED: 09/17/99

SBLKDI	MB1	W	99LE1150	N/A	09/21/99	10/04/99
SBLKDI	MB1 BS	W	99LE1150	N/A	09/21/99	10/04/99

RECRA LabNet Use	Only
99091126	

Custody Transfer Record/Lab Work Request Page 1 of 1 FIELD PERSONNEL: COMPLETE ONLY SHADED AREAS B wetchem

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client Tru Hanford 1399-085						Refrige	erator#	· ,	١	w			<u> </u>	6	6	6	6			*``
Est. Final Proj. Sampling Date						#/Type	Container	Liquid	3v	296	 		4-4	95	-17	4	15			
Project # 18985-001-001-9999-00								Solid	 _							<u> </u>				
Project Contact	Phone	e#				Volum	3	Liquid	Word	1	ļ		 	14	14	11	11-			
RECRA Project	Manag	ger <i>OJ</i>	· 					Solid	ļ	ļ	<u> </u>				_ ZNAC		<u>-</u>			
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Date Rec'd	\J-C	79	Date Due		117/99	ANALY		-	VOA	ORGANIC Sest A					Sufide	AC Prior S	500 505			
MATRIX	T				Matrix	-	TT			2 3	-	R	ECRA L				1			
CODES:	Lab ID	Client 10/De		ription	QC Chose (V)	en Matrix	Date Collected	Time Collected	02024H	\$ 45.55 \$ 45.55				meto	15.2	243O	1 Usus			
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saf 4	130	19-085				met O	2. = O.S.	Bajo	ed,	Cr.Pb,	Se, C	ig, Ci	v	1) Shi Hand	pped _ L Delivered .		1) F	resent or kage	n Oute	r
9/23/99-	INH.	3Nackded	10001	per clu	int coc.		3. My	$V_{1}Zn$	_Be					Airbill	#_*			Inbroken kage (Y)		
	C		-		_9	ang (1)	4- ICCL	10F	-C, 1	CNOZ,	(CNO	3. ICPO	4,	2) Am	bient or Cr	_		resent pa	g Samp	ple
Paf # 1399-085 9/23/99- INH3Nackded to ool per claint cac. COMPOSITE WASTE						5. <u>1056</u>	4.11	2 H $^{\perp}$	NH3N				3) Red Condit	eived in G			(X	or I	N	
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CHAIN OF CUSTODY / REQUEST FOR ANALYSIS RECORD

REFERENCE DOCUMENT NO.: 52-99-0918 CINCINNATI, OH 45253-8704 ORIGINAL MEF NO: FOR SAMPLE RELATED PROBLEMS
ACS CONTACT / PHONE: FULLY REQUIRED REPORT DATE / LAB TAT: SAMPLE SHIPPER (Print): Housen EM 3267 AMPLE NUMBER ANALYSES REQUESTED COLLECTION CONTAINER PRESERVATION SAMPLE MATRIX If more space is required, use FACTS ID **CUSTOMER ID / SAMPLE POINT** DATE TIME 20036161712483-6B-L 200361613 3A4A-SUB-TB5 To sample telementhe TAL I = Jotal volatiles * All TRIP Blanks have RELINQUISHED RECEIVED ITEM / REASON RELINQUISHED BY (Signature) / AFFILIATION BADGE NO. DATE RECEIVED BY (Signature) / AFFILIATION BADGE NO. TIME 9 114 144 0430 ON-SITE - RELEASE FILE / OFF-SITE ANALYTICAL LAB - RETURN TO FEMP YELLOW ON-SITE - DISTRIBUTE AS NEEDED / OFF-SITE ANALYTICAL LAB - RETURN TO FEMP BLUE SAMPLING TECH / PROJECT FILE DISTRIBUTION OF COPIES FS-F-3361 (09-21-95) ★ IF "N", THEN COMPLETE APPLICABLE NON-CONFORMATICE PROCEDURE SHADED AREAS ARE TO BE COMPLETED BY THE SAMPLE RECEIVING GROUP.